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~~subscript n is 0 or 1;~~

~~R⁴ is a member selected from the group consisting of H and C₁-C₆ alkyl;~~

R⁵ is a member selected from the group consisting of H, C₃-C₇ cycloalkyl, C₂-C₆ alkenyl, C₂-C₆ alkyne, phenyl substituted with 0-2 R¹⁵; and a C₁-C₆ alkyl substituted with 0-2 R¹⁸, ~~wherein said C₁-C₆ alkyl optionally contains a heteroatom selected from the group consisting of O, S, S(=O), S(=O)₂ and NR¹⁷;~~

~~Y is a member independently selected from the group consisting of a bond and (CR²⁰R²¹)_m-W-(CR²²R²³)_p;~~

~~subscript p is 1 or 2;~~

~~subscript m is 0 or 1;~~

~~W is a member independently selected from the group consisting of a bond, O, S, S(=O), S(=O)₂ and NR¹²;~~

~~X is selected from the group consisting of C(=O), OC(=O), NR²⁴C(=O) and S(=O)₂;~~

each of R⁶, R⁷, R⁸ and ~~R⁹~~ is independently a member selected from the group consisting of H and C₁-C₄ alkyl;

~~Ar is a member selected from the group consisting of phenyl substituted with 0-3 R²⁹; and 5- to 6-membered heteroaryl containing 1 to 4 heteroatoms each independently a member selected from the group consisting of N, O and S, wherein said heteroaryl is substituted with 0-3 R²⁹;~~

each R¹⁰ is independently a member selected from the group consisting of H, C₃-C₇ cycloalkyl, a C₁-C₃ perfluoroalkyl, a C₁-C₄ alkyl substituted with 0-1 R²⁵, and a phenyl substituted with 0-3 R¹⁵;

each R¹¹ is independently a member selected from the group consisting of H, 'BOC, Cbz, C₃-C₈ cycloalkyl, (C₁-C₆ alkyl)-C(=O)-, (C₁-C₆ alkyl)-S(=O)₂- and a C₁-C₆ alkyl;

each of R¹², R¹³ and R¹⁴ is independently a member selected from the group consisting of H and C₁-C₄ alkyl;

each R¹⁵ is independently a member selected from the group consisting of H, OH, F, Cl, Br, I, CN, NO₂, COOR¹³, C(=O)NR¹³R¹⁴, S(=O)₂NR¹³R¹⁴, acetyl, -SCH₃, -S(=O)CH₃, -S(=O)₂CH₃, NR²⁶R²⁷, C₁-C₆ alkoxy, C₁-C₃ perfluoroalkyl, C₁-C₃ perfluoroalkoxy and a C₁-C₆ alkyl;

each R^{16} is independently a member selected from the group consisting of H, OH, COOR^{13} , $\text{C(=O)NR}^{13}\text{R}^{14}$, $\text{S(=O)}_2\text{NR}^{13}\text{R}^{14}$, acetyl, $-\text{SCH}_3$, $-\text{S(=O)CH}_3$, $-\text{S(=O)}_2\text{CH}_3$, $\text{C}_1\text{-C}_6$ alkoxy, $\text{NR}^{26}\text{R}^{27}$, and a phenyl substituted with 0-3 R^{15} ;

~~R^{17} is a member selected from the group consisting of H and $\text{C}_1\text{-C}_4$ alkyl;~~

each R^{18} is independently a member selected from the group consisting of H, OH, F, Cl, CN, NO_2 , C(=O)OR^{30} , $\text{C(=O)NR}^{13}\text{R}^{14}$, $\text{NR}^{11}\text{R}^{12}$, a $\text{C}_1\text{-C}_3$ perfluoroalkyl, a $\text{C}_1\text{-C}_3$ perfluoroalkoxy, a phenyl substituted with 0-3 R^{15} ; and $\text{C}_3\text{-C}_8$ cycloalkyl;

each R^{19} is ~~[[a]]~~ independently a member selected from the group consisting of $\text{C}_1\text{-C}_4$ alkyl, F, Cl, and $\text{C}_1\text{-C}_4$ alkoxy, CF_3 and OCF_3 ;

~~each of R^{20} , R^{21} , R^{22} and R^{23} is independently a member selected from the group consisting of a bond, H, F, OH, $\text{C}_1\text{-C}_4$ alkyl, and $\text{C}_1\text{-C}_3$ alkylhydroxy;~~

~~R^{24} is a member selected from the group consisting of H and $\text{C}_1\text{-C}_4$ alkyl;~~

each R^{25} is independently a member selected from the group consisting of H, $\text{C}_3\text{-C}_7$ cycloalkyl, and a phenyl substituted with 0-3 R^{15} ;

each R^{26} is independently a member selected from the group consisting of H, $\text{C}_1\text{-C}_4$ alkyl, $(\text{C}_1\text{-C}_4 \text{ alkyl})\text{-C(=O)-}$ and $(\text{C}_1\text{-C}_4 \text{ alkyl})\text{-S(=O)}_2\text{-}$;

each R^{27} is independently a member selected from the group consisting of H and $\text{C}_1\text{-C}_4$ alkyl;

each R^{28} is independently a member selected from the group consisting of H, a $\text{C}_1\text{-C}_6$ alkyl, $\text{C}_3\text{-C}_8$ cycloalkyl, a phenyl substituted with 0-3 R^{15} , and a benzyl substituted with 0-2 R^{15} ;

each R^{29} is independently a member selected from the group consisting of H, F, Cl, Br, I, CN, NO_2 , OR^{28} , SR^{28} , S(=O)R^{28} , $\text{S(=O)}_2\text{R}^{28}$, $\text{S(=O)}_2\text{NR}^{13}\text{R}^{14}$, $\text{NR}^{26}\text{R}^{27}$, acetyl, $\text{C(=O)NR}^{13}\text{R}^{14}$, C(=O)OR^{13} , $\text{C}_1\text{-C}_6$ alkyl, OCHF_2 , SCF_3 , OCF_3 , and $-\text{C(=NH)NH}_2$;

~~alternatively, R^{29} and R^9 are taken together to form a 5- to 7-membered fused heterocyclic ring containing 1-2 heteroatom(s) each independently a member selected from the group consisting of N, O and S; wherein said 5 to 7 membered fused heterocyclic ring is substituted with 0-2 R^{19} ;~~

each R^{30} is independently a member selected from the group consisting of H, $\text{C}_3\text{-C}_7$ cycloalkyl, $\text{C}_1\text{-C}_4$ alkyl substituted with 0-1 R^{25} , and a phenyl substituted with 0-3 R^{15} ;

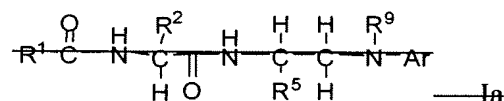
and with the proviso that R^3 , R^4 , R^5 , R^6 , R^7 , and R^8 , ~~and R^9~~ are not all hydrogen.

2-3. (canceled)

4. (previously presented) The compound of claim 1, wherein R¹ is phenyl substituted with 0-3 R^{1a}.

5-6. (canceled)

7. (currently amended) The compound of claim 9 ~~1~~, ~~according to formula Ia:~~



wherein:

R¹ is C₃-C₈ cycloalkyl substituted with 0-2 R^{1b}, wherein said C₃-C₈ cycloalkyl is saturated or unsaturated; and

R² is a member selected from the group consisting of a phenyl substituted with 0-3 R¹⁵, a C₁-C₆ alkyl substituted with 0-2 R^{2a}, and a C₃-C₇ cycloalkyl substituted with 0-2 R¹⁹; and

~~Ar is phenyl substituted with 0-3 R²⁹, or alternatively, R²⁹ and R⁹ are taken together to form a 5- to 7-membered fused heterocyclic ring containing 1-2 heteroatom(s) each independently a member selected from the group consisting of N, O and S; wherein said 5- to 7-membered fused heterocyclic ring is substituted with 0-2 R¹⁹.~~

8. (currently amended) The compound of claim ~~[[7]]~~ 9, wherein:

R² is a member selected from the group consisting of a C₁-C₂ alkyl substituted with 1 R^{2a}, and C₁-C₆ alkyl;

each R^{2a} is independently a member selected from the group consisting of a phenyl substituted with 0-3 R¹⁵, and a C₃-C₈ cycloalkyl substituted with 0-2 R¹⁹;

R⁵ is a member selected from the group consisting of H, C₃-C₇ cycloalkyl; and a C₁-C₆ alkyl substituted with 0-1 R¹⁸; ~~wherein said C₁-C₆ alkyl optionally contains a heteroatom selected from the group consisting of O, S, S(=O), S(=O)₂ and NR¹⁷;~~ and

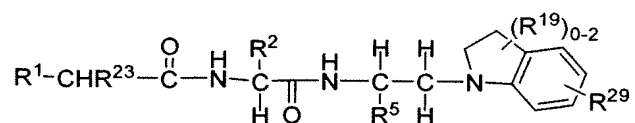
$$\text{R}^1-\text{C}(=\text{O})-\text{NH}-\underset{\text{H}}{\overset{\text{R}^2}{\text{C}}}-\text{C}(=\text{O})-\text{NH}-\underset{\text{R}^5}{\overset{\text{H}}{\text{C}}}-\underset{\text{H}}{\text{C}}-\text{N}(\text{R}^{19})_{0-2}-\text{C}_6\text{H}_3(\text{R}^{29})$$
$$\text{R}^1-\text{CHR}^{23}-\text{C}(=\text{O})-\text{N}(\text{H})-\text{C}(\text{H})(\text{R}^2)-\text{C}(\text{H})(\text{R}^5)=\text{C}(\text{H})(\text{R}^9)-\text{N}(\text{H})-\text{Ar}$$

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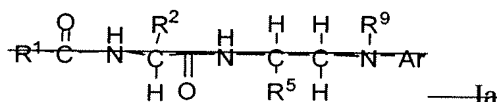
each R^{2a} is independently a member selected from the group consisting of a phenyl substituted with 0-3 R^{15} , and a C_3 - C_8 cycloalkyl substituted with 0-2 R^{19} ; and

R^5 is a member selected from the group consisting of H, C_3 - C_7 cycloalkyl; and a C_1 - C_6 alkyl, ~~wherein said C_1 - C_6 alkyl optionally contains a heteroatom selected from the group consisting of O, S, $S(=O)$, $S(=O)_2$ and NR^{17} .~~

18. (currently amended) The compound of claim [[16]] 1, wherein said compound is of the formula:



19. (currently amended) The compound of claim 9 1, ~~according to formula Ia~~



wherein:

R^1 is C_6 - C_{10} aryl substituted with 0-3 R^{1a} ;

each R^{1a} is independently a member selected from the group consisting of H, C_1 - C_3 perfluoroalkyl, C_3 - C_7 cycloalkyl, F, Cl, Br, CN, NO_2 , OR^{10} , SCH_3 , $S(=O)CH_3$, $S(=O)_2R^{10}$, $NR^{11}R^{12}$, acetyl, $C(=O)OR^{13}$, $C(=O)NR^{13}R^{14}$, $S(=O)_2NR^{13}R^{14}$, phenyl substituted with 0-3 R^{15} ; and a C_1 - C_4 alkyl substituted with 0-2 R^{16} ;

R^2 is a member selected from the group consisting of a phenyl substituted with 0-3 R^{15} ; a C_1 - C_2 alkyl substituted ~~[[with 1 R^{2a}]]~~ with 1 R^{2a} , and a C_3 - C_7 cycloalkyl substituted with 0-2 R^{19} ; and

each R^{2a} is independently a member selected from the group consisting of a C_6 - C_{10} aryl substituted with 0-3 R^{15} ; a C_3 - C_8 cycloalkyl substituted with 0-2 R^{19} ; and a C_7 - C_{11} bicycloalkyl substituted with 0-2 R^{19} ; and

~~Ar is phenyl substituted with 0-3 R^{29} , or alternatively, R^{29} and R^9 are taken together to form a 5- to 7-membered fused heterocyclic ring containing 1-2 heteroatom(s) each~~

~~independently a member selected from the group consisting of N, O and S; wherein said 5- to 7-membered fused heterocyclic ring is substituted with 0-2 R¹⁹.~~

20-26. (canceled)

27. (currently amended) A pharmaceutical composition comprising the compound of Formula I(i) or I(j) in claim 1[: or]] and a pharmaceutically acceptable ~~salt and an~~ excipient.

28. (currently amended) A pharmaceutical composition comprising the compound of claim 38 and a pharmaceutically acceptable excipient.

29-37. (canceled)

38. (currently amended) The compound of claim 1, selected from the group consisting of:

(S)-N-{1-[2-(5-Fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-3-methyl-butyl}-3-methyl-benzamide;

N-(S)-{2-cyclohexyl-1-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-4-phenoxy-benzamide;

(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-[2-(4-methoxy-phenyl)-acetylamino]-propionamide;

(S)-N-{1-[2-(5-Chloro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-3-methyl-butyl}-3-methyl-benzamide;

(S)-N-{3-Cyclohexyl-1-[2-(7-methoxy-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;

(S)-N-{3-Cyclohexyl-1-[2-(6-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;

(S)-N-{3-Cyclohexyl-1-[2-(7-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;

(S)-N-{3-Cyclohexyl-1-[2-(5-cyano-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;

Cyclopropanecarboxylic acid (S)-{2-cyclohexyl-1-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-amide;

(S)-N-{3-Cyclohexyl-1-[2-(4-methoxy-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;

(S)-N-{3-Cyclohexyl-1-[2-(5-methoxy-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;

(S)-N-{3-Cyclohexyl-1-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;

(S)-N-{3-Cyclohexyl-1-[2-(5-benzyloxy-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;

~~N-{1-(S)-[2-(4-Methoxy-phenylamino)-propylcarbamoyl]-3-methyl-butyl}-3-methyl-benzamide;~~

~~N-{1-(S)-[2-(4-Methoxy-phenylamino)-1-methyl-ethylcarbamoyl]-3-methyl-butyl}-3-methyl-benzamide;~~

~~N-{1-(S)-[2-(4-Methoxy-phenylamino)-1-(S)-methyl-ethylcarbamoyl]-3-methyl-butyl}-3-methyl-benzamide;~~

~~N-{1-(S)-[2-(4-Methoxy-phenylamino)-1-(R)-methyl-ethylcarbamoyl]-3-methyl-butyl}-3-methyl-benzamide;~~

~~N-{2-Cyclohexyl-(1S)-[2-(4-methoxy-phenylamino)-(1R)-methyl-ethylcarbamoyl]-ethyl}-3-methoxy-benzamide;~~

~~N-{(1S)-[2-(5-Fluoro-2,3-dihydro-indol-1-yl)-1,1-dimethyl-ethylcarbamoyl]-2-phenyl-ethyl}-3-methyl-benzamide;~~

~~N-{1-(S)-[1-(R)-Benzyloxymethyl-2-(4-methoxy-phenylamino)-ethylcarbamoyl]-3-methyl-butyl}-3-methyl-benzamide;~~

~~N-(S)-{[1-(R)-Benzyloxymethyl-2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-phenyl-methyl}-3-methoxy-benzamide;~~

~~N-[1-(S)-[1-(R)-Benzyloxymethyl-2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-2-(4-fluoro-phenyl)-ethyl]-3-methoxy-benzamide;~~

~~N-{1-(S)-[(2-Benzyloxy-1-(R)-(5-fluoro-2,3-dihydro-indol-1-yl)methyl)-ethylcarbamoyl]-3-cyclohexyl-propyl}-3-methoxy-benzamide;~~

~~N-{3-Cyclohexyl-1-(S)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-(R)-hydroxymethyl-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;~~

N-{3-Cyclohexyl-1-(*R*)-[(*S*)-2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-(*R*)-hydroxymethyl-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;

(*S,S*)-5-(5-Fluoro-2,3-dihydro-indol-1-yl)-4-[4-methyl-2-(3-methyl-benzoylamino)-pentanoylamino]-pentanoic acid benzyl ester;

(*S,S*)-5-(5-Fluoro-2,3-dihydro-indol-1-yl)-4-[4-methyl-2-(3-methyl-benzoylamino)-pentanoylamino]-pentanoic acid;

(*S,S*)-*N*-{1-[3-Carbamoyl-1-(5-fluoro-2,3-dihydro-indol-1-yl)methyl]-propylcarbamoyl]-3-methyl-butyl}-3-methyl-benzamide;

(*S,S*)-~~*N*-{1-[1-(5-Fluoro-2,3-dihydro-indol-1-yl)methyl)-3-ureido-propylcarbamoyl]-3-methyl-butyl}-3-methyl-benzamide;~~

(*S,S*)-3-[4-Cyclohexyl-2-(3-methoxy-benzoylamino)-butyrylamino]-4-(5-fluoro-2,3-dihydro-indol-1-yl)-butyric acid benzyl ester;

(*S,S*)-3-[4-Cyclohexyl-2-(3-methoxy-benzoylamino)-butyrylamino]-4-(5-fluoro-2,3-dihydro-indol-1-yl)-butyric acid;

(*S,S*)-*N*-{1-[1-Benzyl-2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-3-cyclohexyl-propyl}-3-methoxy-benzamide;

(*S,S*)-*N*-{3-Cyclohexyl-1-[1-(5-fluoro-2,3-dihydro-indol-1-yl)methyl)-3-methyl-butylcarbamoyl]-propyl}-3-methoxy-benzamide;

(*S,S*)-*N*-{3-Cyclohexyl-1-[1-(5-fluoro-2,3-dihydro-indol-1-yl)methyl)-2-methyl-propylcarbamoyl]-propyl}-3-methoxy-benzamide;

(*S,S*)-*N*-{3-Cyclohexyl-1-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-phenyl-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;

~~*N*-{1-(*S*)-[2-(*R*)-Benzyloxy-1-(*R*)-(5-fluoro-2,3-dihydro-indol-1-yl)methyl)-propylcarbamoyl]-3-cyclohexyl-propyl}-3-methoxy-benzamide;~~

~~*N*-{1-(*R*)-[1-(*R*)-Benzylsulfanylmethyl-2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-3-cyclohexyl-propyl}-3-methoxy-benzamide;~~

(*S,S*)-{5-[4-Cyclohexyl-2-(3-methoxy-benzoylamino)-butyrylamino]-6-(5-fluoro-2,3-dihydro-indol-1-yl)-hexyl}-carbamic acid benzyl ester;

N-{2-cyclohexyl-1-(*S*)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-2-(2-fluoro-biphenyl-4-yl)-propionamide;

N-{2-cyclohexyl-1-(*S*)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-2-*p*-tolyl-propionamide;

N-{2-cyclohexyl-1-(S)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-
2-o-tolyl-propionamide;

N-{2-cyclohexyl-1-(S)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-
2-(4-fluoro-phenyl)-propionamide;

2-(4-Chloro-phenyl)-N-{2-cyclohexyl-1-(S)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-
ethylcarbamoyl]-ethyl}-propionamide;

N-{2-cyclohexyl-1-(S)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-
2-(R)-phenyl-propionamide;

N-(S)-{2-cyclohexyl-1-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-
3-methyl-benzamide;

N-(S)-{2-cyclohexyl-1-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-
4-(methanesulfonylamino-methyl)-benzamide;

N-(S)-{2-cyclohexyl-1-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-
3-methanesulfonyl-benzamide;

N-(S)-{2-cyclohexyl-1-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-
4-methanesulfonylamino-benzamide;

N-{2-cyclohexyl-1-(S)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-
2-(4-hydroxy-phenyl)-propionamide;

4-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-(S)-(2-(R)-phenyl-
propionylamino)-butyramide;

N-{2-cyclohexyl-1-(S)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-
2-(R)-phenyl-butyramide;

~~N-{1-(S)-[1-(R)-Benzyloxymethyl-2-(5-fluoro-2,3-dihydro-indol-1-yl)-
ethylcarbamoyl]-2-cyclohexyl-ethyl}-3-methoxy-benzamide;~~

N-{2-Cyclohexyl-1-(S)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-(R)-hydroxymethyl-
ethylcarbamoyl]-ethyl}-3-methoxy-benzamide;

~~N-{1-(S)-[1-(R)-Benzyloxymethyl-2-(5-fluoro-2,3-dihydro-indol-1-yl)-
ethylcarbamoyl]-3,3-dimethyl-butyl}-3-methoxy-benzamide;~~

N-{1-(S)-[2-(5-Fluoro-2,3-dihydro-indol-1-yl)-1-(R)-hydroxymethyl-
ethylcarbamoyl]-3,3-dimethyl-butyl}-3-methoxy-benzamide;

~~3-(S)-(2-(S)-Benzyloxy-carbonylamino-4,4-dimethyl-pentanoylamino)-4-(5-fluoro-
2,3-dihydro-indol-1-yl)-butyric acid tert-butyl ester;~~

~~3-(S)-(2-(S)-Benzyloxycarbonylamino-4,4-dimethyl-pentanoylamino)-4-(5-fluoro-2,3-dihydro-indol-1-yl)-butyric acid;~~

~~4-(5-Fluoro-2,3-dihydro-indol-1-yl)-3-(S)-[2-(S)-(3-methoxy-benzoylamino)-4,4-dimethyl-pentanoylamino]-butyric acid tert-butyl ester;~~

~~3-(S)-[3-Cyclohexyl-2-(S)-(3-methoxy-benzoylamino)-propionylamino]-4-(5-fluoro-2,3-dihydro-indol-1-yl)-butyric acid benzyl ester;~~

~~3-(S)-[3-Cyclohexyl-2-(S)-(3-methoxy-benzoylamino)-propionylamino]-4-(5-fluoro-2,3-dihydro-indol-1-yl)-butyric acid;~~

~~4-(5-Fluoro-2,3-dihydro-indol-1-yl)-3-(S)-[2-(S)-(3-methoxy-benzoylamino)-4,4-dimethyl-pentanoylamino]-butyric acid ethyl ester;~~

~~N-{1-(S)-[2-Cyano-1-(S)-(5-fluoro-2,3-dihydro-indol-1-ylmethyl)-ethylcarbamoyl]-3,3-dimethyl-butyl}-3-methoxy-benzamide;~~

~~N-{1-(S)-[5-Amino-1-(S)-(5-fluoro-2,3-dihydro-indol-1-ylmethyl)-pentylcarbamoyl]-3-cyclohexyl-propyl}-3-methoxy-benzamide;~~

~~3-(S)-(2-(S)-Benzyloxycarbonylamino-3-cyclohexyl-propionylamino)-4-(5-fluoro-2,3-dihydro-indol-1-yl)-butyric acid benzyl ester;~~

~~1-(S)-[1-(R)-Benzyloxymethyl-2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-2-cyclohexyl-ethyl}-carbamie acid benzyl ester;~~

~~N-{3-Cyclohexyl-1-(S)-[2-(3,5-dimethoxy-benzyloxy)-1-(R)-(5-fluoro-2,3-dihydro-indol-1-ylmethyl)-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;~~

~~4-{2-(R)-[4-Cyclohexyl-2-(S)-(3-methoxy-benzoylamino)-butyrylamino]-3-(5-fluoro-2,3-dihydro-indol-1-yl)-propoxymethyl}-benzoic acid methyl ester;~~

~~(S,S)-N-{3-Cyclohexyl-1-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-(4-hydroxy-benzyl)-ethylcarbamoyl]-propyl}-3-methoxy-benzamide;~~

~~{2-Cyclohexyl-1-(S)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-(S)-methyl-ethylcarbamoyl]-ethyl}-carbamie acid benzyl ester;~~

~~4-Benzyloxy-N-(R,S)-{[2-(4-amidinophenylamino)-1-(S)-methyl-ethylcarbamoyl]-(2,4-dichloro-phenyl)-methyl}-benzamide;~~

~~{1-(S)-[2-(5-Fluoro-2,3-dihydro-indol-1-yl)-1-(S)-methyl-ethylcarbamoyl]-3,3-dimethyl-butyl}-carbamie acid benzyl ester;~~

~~Cyclopropanecarboxylic acid {1-(S)-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-(S)-methyl-ethylcarbamoyl]-3,3-dimethyl-butyl}-amide;~~

~~(S,S)-2-(3-Chloro-benzenesulfonylamino)-3-cyclohexyl-N-[1-methyl-2-(4-trifluoromethoxy-phenylamino)-ethyl]-propionamide;~~

~~(S,S)-3-Cyclohexyl-N-[1-methyl-2-(4-trifluoromethoxy-phenylamino)-ethyl]-2-(3-trifluoromethoxy-benzenesulfonylamino)-propionamide;~~

N-((S)-(2-(5-fluoroindolin-1-yl)ethylcarbamoyl)(cyclohexyl methyl)-3-methylbenzamide;

N-((S)-1-(2-(5-fluoroindolin-1-yl)ethylcarbamoyl)-2-(2-chlorophenyl)ethyl)-3-methylbenzamide;

N-((S)-1-(2-(5-fluoroindolin-1-yl)ethylcarbamoyl)-2-(3-chlorophenyl)ethyl)-3-methylbenzamide;

N-((S)-1-(2-(5-fluoroindolin-1-yl)ethylcarbamoyl)-2-(4-chlorophenyl)ethyl)-3-methylbenzamide;

(S)-N-{2-Cyclopentyl-1-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethylcarbamoyl]-ethyl}-3-methyl-benzamide;

N-((S)-1-(2-(5-fluoroindolin-1-yl)ethylcarbamoyl)-3,3-dimethylbutyl)-3-methylbenzamide;

N-((S)-1-(2-(5-fluoroindolin-1-yl)ethylcarbamoyl)-3-cyclohexylpropyl)-3-methylbenzamide;

N-((S)-1-(2-(5-fluoroindolin-1-yl)ethylcarbamoyl)-2-phenylethyl)-3-methylbenzamide;

N-(R,S)-((3-(5-fluoroindolin-1-yl)-1-hydroxypropan-2-(R)-ylcarbamoyl)(2,4-dichlorophenyl)methyl)-3,4-difluorobenzamide;

~~N-(S)-((3-(benzyloxy)-1-(5-fluoroindolin-1-yl)propan-2-(R)-ylcarbamoyl)(2,4-dichlorophenyl)methyl)-3,4-difluorobenzamide;~~

(R,S)-N-((2-(5-fluoroindolin-1-yl)ethylcarbamoyl)(2,4-dichlorophenyl)methyl)-3-methylbenzamide;

(S,S)-N-((3-(5-fluoroindolin-1-yl)-1-hydroxypropan-2-ylcarbamoyl)(2,4-dichlorophenyl)methyl)-3,4-difluorobenzamide;

(S,S)-4-(5-Fluoro-2,3-dihydro-indol-1-yl)-3-[2-(3-methoxy-benzoylamino)-4,4-dimethyl-pentanoylamino]-butyric acid; and

~~(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-(5-isoxazol-3-yl-thiophene-2-sulfonylamino)-propionamide;~~

~~(S)-2-(3-Biphenyl-4-yl-ureido)-3-cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-propionamide;~~

~~(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-(4-phenoxy-benzenesulfonylamino)-propionamide;~~

~~(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-(naphthalene-1-sulfonylamino)-propionamide;~~

~~(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-(4-trifluoromethyl-benzenesulfonylamino)-propionamide;~~

~~(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-(4-trifluoromethoxy-benzenesulfonylamino)-propionamide;~~

~~(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-[4-(4-fluorophenoxy)-benzenesulfonylamino]-propionamide;~~

~~(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-(4'-methoxybiphenyl-4-sulfonylamino)-propionamide;~~

~~(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-(4-methoxy-benzenesulfonylamino)-propionamide;~~

~~(S)-3-Cyclohexyl-2-(4-difluoromethoxy-benzenesulfonylamino)-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-propionamide;~~

~~(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-phenylmethanesulfonylamino-propionamide;~~

~~(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-(toluene-3-sulfonylamino)-propionamide;~~

~~(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-[4-(4-methoxyphenoxy)-benzenesulfonylamino]-propionamide;~~

~~(S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-ethyl]-2-(3-methoxy-benzenesulfonylamino)-propionamide;~~

~~(S,S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-methyl-ethyl]-2-(toluene-3-sulfonylamino)-propionamide;~~

~~(S,S)-3-[4,4-Dimethyl-2-(toluene-3-sulfonylamino)-pentanoylamino]-4-(5-fluoro-2,3-dihydro-indol-1-yl)-butyric acid tert-butyl ester;~~

~~(S,S)-3-Cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-methyl-ethyl]-2-(3-trifluoromethoxy-benzenesulfonylamino)-propionamide;~~

~~(S,S)-2-(3-Chloro-benzenesulfonylamino)-3-cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-methyl-ethyl]-propionamide;~~

~~(S,S)-N-{3-Cyclohexyl-1-[1-(5-fluoro-2,3-dihydro-indol-1-ylmethyl)-3-hydroxy-propylcarbamoyl]-propyl}-3-methoxy-benzamide;~~

~~(S,S)-3-[4,4-Dimethyl-2-(toluene-3-sulfonylamino)-pentanoylamino]-4-(5-fluoro-2,3-dihydro-indol-1-yl)-butyric acid;~~

~~(S,S)-2-Benzenesulfonylamino-3-cyclohexyl-N-[2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-methyl-ethyl]-propionamide; and~~

~~(S,S)-4,4-Dimethyl-2-(toluene-3-sulfonylamino)-pentanoic acid [2-(5-fluoro-2,3-dihydro-indol-1-yl)-1-methyl-ethyl]-amide.~~